Ref #	Hits	Search Query	DBs	Default Operat or	Plura Is	Time Stamp
L1	460523	(alumina or "activated alumina" or aluminate or aluminate or aluminate or aluminate or aluminate or bauxite or "aluminium oxide" or corundum or "IAI2O3!" or "\$AI2O3\$" or boehmite or diaspore or aluminasol or aluminasol\$)	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 12:48

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		EAST Sea	rch Histor	ry		
7/11/24	2455	L1 AND (Tetracycline or tetracyclin or abramycin or abramycin or abramycin or abramycin or achromycine or achromycine or achromycine or achromycine or acromicina or actisite or agromicina or atsite or agromicina or atsite or ambramicina or ambramicine or ambramycine or ambramycine or ambramycine or "apo-tetra" or "apo-tetra" or "apo-tetra" or "apo-tetra" or "bio tetra" or biocycline or biocycline or biocycline or biocycline or biocycline or calocicina or calocicina or cyclomycin or cyclomycin or cyclomycin or deschlorobiomycine or deschlorobiomycine or dispatetrine or economycine or economycine or enterocycline or enterocycline or helvecycline or helvecycline or hexacycline or hexacycline or lidy or l	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 11:59
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		EAST Sear		. ,			
L3	1177	L1 AND ((Albofungin) OR (Antimycin or Antimycine)	US-PGPU B;	OR	ON	2006/07/11 12:00	
		OR (Baicalein or Baicalin	USPAT;				
		or Trihydroxyflavone or	EPO;		ł		
		Baikalein or Noroxylin) OR	JPO; DERWEN				
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		Ibidomide) OR					l
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	ł	Zestocarp) OR (Osalmid or		1			l
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		Bilene or Bilocol or Dribazil		1			l
		or Driol or "Driol-Labaz" or		1	1		
		"Driol Labaz" or DriolLabaz		1	1		
		or Enidran or Fumispore		1	1		
		or Jestmin or Kanochol or "!Li Dan Feng!" or		1	1		
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1		e or Hydroxysalicylanilide		1		l	
		or Salicoylaminophenol or					
		Salmidochol or Saryuurin				1	
	L	or Yoshicol))	L	<u></u>		L	J

				<u> </u>			
	1890	L1 AND ((Salacetamide or Acetsalicy/amide or Acetsalicy/amide or Actylamide or Actylamide or Arthrisin or Ethrisin or Labazyl or Nacemide or Rixamone or Salicyl) OR (Salicy/amide or Acket or Afko-Sal" or "Afko Sal" or Afko-Sal" or "Afko Sal" or Amid-Sal" or Algamon or Algiamida or Algamon or Algiamida or Algamon or Algamida or Amid-Sal" or Or "Amid-Sal" or Amid-Sal" or Amid-Sal or Amid-Sal or Amid-Sal or Or Carboxamidophenol or Carboxamidophenol or Cetamide or Cidal or Cymidon or Dropsprin or Liquiprin or Morsarinas or Novecyl or Hydroxybenzamide or Oramid or Panithal or Raspberin or Salamid or Salimid or Salicylic acid amide" or Salicylic acid amide" or Salicylarilide or Ansadol or Aseptolan or Hyanliid or Asetylamide or phenylsalicylamide or Salicylarilido or Disalde o	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 12:00 Page 4 palized Alumina Particle	
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L5	7571						İ
_		L1 AND ((AlizarIn or	US-PGPU	OR	ON	2006/07/11	
	, ,,,,	Alizarine or	В;	•	0.,	12:39	
		Anthraquinonediol or	USPAT;				
		dihydroxyanthracenedione	EPO;				
		or	JPO;				
		dihydroxyanthraquinone	DERWEN				
		or "Deep Crimson Madder"	Τ				
		or "Eljon Madder" or					
		"Metachrome Red" or					
		"Mordant Red" or "Turkey					
		Red") OR (Alkannin or					
		Alkannine or Alkanet or					
		"Alkanet extract" or					
		Alkanna or "Alkanna red"					
		or Alkannin or Anchusa or					
		"Anchusa acid" or					
		Anchusin or "!Natural Red					
		20!") OR (Anthragaliol or					
		Antragallol or					
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		or "Alizarin Brown" or					
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		"Anthracene Brown" or					
		"Antracromo Brown" or					
		"!Mordant Brown 42!" or					
		"Chrome Fast Brown" or					
		"Mitsul Anthracene					
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		or Anthrarufine or	ł	l			
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		or		1			
		dihydroxyanthraquinone)		1			
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		Altan or Antrapurol or		1			
)	Chrysazin or Diaquone or					l
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		OF dihydroxyanthraguinone		1	1		
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		Dionone or Dorbane or		1			
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		or Laxipur or Laxipurin or					
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L6	11214	L2 OR L3 OR L4 OR L5	US-PGPU 8; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 12:43
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L12	759	L9 AND ((alumina or "activated alumina" or aluminate or aluminated or "aluminum oxide" or bauxite or "aluminum oxide" or bauxite or "aluminum oxide" or bauxite or "aluminum oxide" or boehmite or diaspore or aluminasol or aluminasols) NEAR30 (micrometer or "micrometer" or "micrometers or "micrometers" or "micrometers" or "micrometers" or "micrometers" or micronize or "nanometers" or "nanometers" or "nanometers" or "nanometers" or "nanometers" or "nanometers" or "nanometers" or "nanometers" or "picometers" or "picometers or "picometers" or "picometers or "picometers" or "picometers" or "pico or angstromorometers or aluminated or angstromorometers)	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 12:52
L13	4243	L10 OR L11 OR L12	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 12:53

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EAST Search History

L28	128	MacDonald-John-Gavin.IN.	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 21:10
L29	22	Lye-Jason.IN.	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 21:11
L30	8711	Kimberly-Clark.AS.	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 21:11
L31	9437	L26 OR L27 OR L28 OR L29 OR L30	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 21:12
L32	7966	L31 AND @AD<="20021220"	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 21:14
L33	10	L32 AND L13	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/11 21:14

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Ref #	Hits	Search Query	DBs	Default Operat or	Plura Is	Time Stamp
L1	102	(transdermal) NEAR5 (antimicrobial)	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/12 09:49

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L6	13	L5 AND @AD<="20021220"	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/12 09:57
L7	2	L6 AND alumina	US-PGPU B; USPAT; EPO; JPO; DERWEN T	OR	ON	2006/07/12 09:57

7/12/2006 10:16:21 AM Page

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2-wapnthacenecarboxamide, 4-(dimethylamino)-1,4,4a,5,5a,6,11,12a-octahydro-3,6,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo-, (4S,4aS,5aS,6S,12aS)-
                                                                                                               2-Naphthacenecarboxamide, 4-(dimethylamino)-1,4,4a,5,5a,6,11,12a-octahydro-3,6,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo-, [45-
(4α,4aα,5aα,6β,12aα)}-
                                                                                                                                           2-Naphthacenecarboxamide, 4-(dimethylamino)-1,4,4a,5,5a,6,11,12a-octahydro-3,6,10,12,12a-pentahydroxy-6-methyl-1,11-dioxo- (7CI, 8CI)
                                                COPYRIGHT 2006 ACS on STN
                                                                                                                                                                                            Achromycin
Achromycin (naphthacene derivative)
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1 TETRACYCLINE/CN
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C22 H24 NZ 08
C22 H24 NZ 08
TO Files: ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BELLSTEIN*, BIOSIS,
BIOTECHNO, CA, CABA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMATS, CHEMIST,
CIN, CSCHEM, CSNB, DDFU, DRUGU, EMBASE, GMELIN*, HSDB*, IFICDB, IFICDB, IFICDB, IFICDB, IFICDB, IFICDB, IPICOSCARCH, IPA, MEDLINE, MRCK*, MSDS-OHS, NAPRALERT, PHAR,
PIRM, PROMT, PS, RTECS*, SPECINFO, TOXCENTER, USAN, USPATZ, USPATFULL,
VETU, VTB
CN Retet
CN Robitet Robicaps
CN Tetracycline
ADDITIONAL NAMES NOT AVAILABLE IN THIS FORMAT - Use FCN, FIDE, or ALL for
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                         (File contains numerically searchable property data)
r Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)
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823 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE
16149 REFERENCES IN FILE CAPLUS (1907 TO DATE)
I REFERENCES IN FILE CAPLUS (1907 TO 1967)
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    **PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                Absolute stereochemistry. Rotation (-).
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                 -> S Albofungin/CN
12 1 ALBOFUNGIN/CN
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                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                      퓽
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CN Albofung
CN Antibiot
DR 37184-49
MF C27 H24
LC STN File
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                        => D L2
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19 REFERENCES IN FILE CA (1907 TO DATE) 19 REFERENCES IN FILE CAPLUS (1907 TO DATE)

-> S Antimycin Al/CN L3 1 ANTIMYCIN Al/CN -> D L3 1

L3 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN
RN 642-15-9 REGISTRY
ED Entered STNR: 16 Nov 1984
CN Butanoic acid, 2(or 3)-methyl-, (2R,3S,6S,7R,8R)-3-[[3-(formylamino)-2-hydranoic acid, 2(or 3)-methyl-, (2R,3S,6S,7R,8R)-3-[[3-(formylamino)-2-hydranoic acid, 2(or 3)-methyl-, 6-dimethyl-4,9-dioxo-1,5-dioxonan-7-yl ester (9C1) (CA INDEX NAMES:
CN 1,5-Doxonano, butanoic acid deriv.
CN Antimycin A1 (6C1, 7C1)
CN Antimycin A1 (6C1, 7C1)
CN Antimycin A1 (6C1, 7C1)
CN Antimycin A1 (6C1, 7C1)
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CN Antimycin A1 (6C1, 7C1)
CN Antimycin

D1-Me

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

37 REFERENCES IN FILE CA (1907 TO DATE)
5 REFERENCES TO NON-SPECIFIC DENYATIVES IN FILE CA 37 REFERENCES IN FILE CAPLUS (1907 TO DATE)
11 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> S Baicalein/CN L4 1 BAICALEIN/CN

-> D L4 1

ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 491-67-8 REGISTRY Entered STN: 16 Nov 1984 4H-1-Benzopyran-4-one, 5,6,7-trihydroxy-2-phenyl- (9CI) (CA INDEX NAME)

Baicalein (6CI) Flavone, 5,6,7-trihydroxy- (7CI, 8CI) 5, 6, 7-Trihydroxyflavone

LLA ANSWER 1

RN 491-67-6

ED F10-67-6

CN 4H-1-BEC

CN F12VONE,

CN F12VONE,

CN BAIKALEI

CN BAIKALEI

CN BAIKALEI

CN NOCOXYII

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TW Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BELLSTEIN*, BIOSIS, BIOTECHNO, CA., CAOLD, CAPLUS, CASREACT, CHEMCATS, CSCHEM, DDFU, DRUGU, EMBASE, IPA, MEDLINE, MRCK*, NAPRALERT, PHAR, PROUSDDR, RTECS*, TOXCENTER, USPAT2, USPATPLL.

(*File contains numerically searchable property data) Baikalein Noroxylin NSC 661431 3D CONCORD C15 H10 05 COM STN Files: 1

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

922 REFERENCES IN FILE CA (1907 TO DATE) 9 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

929 REFERENCES IN FILE CAPLUS (1907 TO DATE) 12 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> S Baicalin Hydrate/CN L5 0 BAICALIN HYDRATE/CN

S Baicalin/CN 1 BAICALIN/CN

=> D L6 1

CERE

ANSWER I OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 21967-41-9 REGISTRY Thereed STN: 16 Nov 1984 B-D-Glucopyranosidutonic acid, 5,6-dihydroxy-4-oxo-2-phenyl-4H-1-benzopyrano-7-yl (9CI) (CA INDEX NAME)

Baicalin (6CI, 7CI, 8CI)

OTHER CA INDEA CN Batcalin OTHER NAMES: CN Batcalein CN Batcalein CN Batcalein FS STEREOSEI MF C21 H18 (

Marcalein 7-glucuronide
Baicalein 7-G-\$\beta-D-glucuronide
Baicalein 7-O-\$\beta-D-glucuronide
Baicalein 7-O-\$\beta-D-glucuronide
STEREOSEARCH
27462-75-5, 31564-28-0, 100647-26-5
C21 H18 011
COM
FILES: ADISINSIGHT, AGRICOLA, ANABSTR, BELLSTEIN*, BIOSIS,
STN FILES: ADISINSIGHT, AGRICOLA, ANABSTR, RELSTEIN*, DEDU,
BUOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMCATS, CSCHEM, DEDU,
BUNGE, IPA, MEDLINE, NAPRALERT, PROUSDDR, RTECS*, TOXCENTER,
USPATZ, USPATFULL
(*File contains numerically searchable property data)

Absolute stereochemistry

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

993 REFERENCES IN FILE CA (1907 TO DATE)
23 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1006 REFERENCES IN FILE CAPLUS (1907 TO DATE)
9 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> S Carubicin/CN 1 CARUBICIN/CN

-> D L7 1

ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 50935-04-1 REGISTRY Entered STN: 16 Nov 1984 228

CN 5,12-Naphthacenedione, 8-acetyl-10-[(3-amino-2,3,6-trideoxy-α-1-1yxo-hexcopyranosyl)oxyl-7,8,9,10-tetrahydro-1,6,8,11-tetrahydroxy-, (85,105)-(9CI) (CA INDEX NAME) OTHER CA INDEX NAMES:

Z

5,12-Naphthacenedione, 8-acetyl-10-[(3-amino-2,3,6-trideoxy-α-L-1yxo-hexopyranosyl)oxyl-7,8,9,10-tetrahydro-1,6,8,11-tetrahydroxy-, (8S-cis)-

Antibiotic R 588A Carminomicin I Carminomycin I

OTHER NAMES:
CN Antibiot
CN Carminot
CN Carminot
CN Carminot
CN Carminot
CN Carubict
FS STEREOSE
MF C26 H27
LC COM

Carubicin.
Carubicin.
STEREOSEARCH
CZ6 H27 N O10
CCM
STN Files: AGRICOLA, ANABSTR, BELLSTEIN*, BIOSIS, BIOTECHNO, CA, CAPLUS,
CASREACT, DDFU, DRUGU, EMBASE, IFICOB, IFIPAT, IFIUDB, IPA, MRCK*,
NAPRALERT, PHAR, RIECS*, TOXCENTER, USAN, USPAT2, USPATFUL
(*File contains numerically searchable property data)

Absolute stereochemistry

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

225 REFERENCES IN FILE CA (1907 TO DATE)
27 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
228 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> S Closantel/CN L8 1 CLOSANTEL/CN

=> D L8 1

ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 57808-65-8 REGISTRY CARLES OF STN: 16 Nov 1994 Benzamide, N-[5-chloro-4-[(4-chlorophenyl)cyanomethyl]-2-hydroxy-3,5-diiodo- (9CI) (CA INDEX NAME) 3 Z B 3

OTHER NAMES

Closantel NSC 335306 Zycloz 3D CONCORD

2252

C 3

CC2 H14 C12 12 N2 O2
COM
TOWN
STN Files: AGRICOLA, ANABSTR, AQUIRE, BIOSIS, BIOTECHNO, CA, CABA,
CRAPLUS, CRNB, CHEMCATS, CHEMLIST, CSCHEM, DDFU, DRUGU, ENBASE, IFICEDB,
IFIPRIT, IFIUDB, IFP, MEDLINE, MRCK*, PROMT, RIECS*, TOXCENTER, USAN,
USPATFULL, VETU

(*File contains numerically searchable property data)
Other Sources: EINECS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

144 REFERENCES IN FILE CA (1907 TO DATE) 8 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA 144 REFERENCES IN FILE CAPLUS (1907 TO DATE)

S £ 5

Daunorubicin/CN
1 DAUNORUBICIN/CN

D L9 1 î ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 20830-81-3 REGISTRY

Entered STN: 16.00v 1984 5,12-Naphthacendione, 8-acetyl-10-[(3-amino-2,3,6-trideoxy-α-L-lyxo-hexopyranosyl)oxy]-7,8,9,10-tetrahydro-6,8,11-trihydroxy-1-methoxy-,(85,10S)- (9E1) (CA INDEX NAME) CERE

5,12-Naphthacenedione, 8-acetyl-10-[(3-amino-2,3,6-trideoxy- α -L-1yxo-hexopyranosyl)oxy]-7,8,9,10-tetrahydro-6,8,11-trihydroxy-1-methoxy-, (8S-cis)-OTHER CN 5

Daunomycin (8CI) NAMES:

Acetyladriamycin (+)-Daunomycin

Daunoblastina Cerubidin

Daunomycine Daunorubicin Daunorubicine DaunoXome

Leukaemomycin C NSC 82151 NSC 83142

Rubidomycin

Rubomycin C STEREOSEARCH 11006-64-5, 11048-29-6, 1407-15-4, 23942-76-9, 149541-57-1, 27576-81-4, 28020-80-6 C27 H29 N 010

COM STN Files: 53

STU Files: ADISINSIGHT, ADISNEMS, AGRICOLA, ANABSTR, BELLSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASEACT, CBNB, CHEMCATS, CHEMINFORMEX, CHEMLIST, CIN, CSCHEM, CSNF, DDGU, BRABSE, HSDB., IFIDEB, IFIPAT, IFIUDB, IMSCOSEARCH, ITEA, MEDLINE, MACK+, MEDASCH, MSDB-, IRICORE, IFIPAT, PIRA, PROMT, PS, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, USPAT2, USPATFULL (*File contains numerically searchable property data) Other Sources: BINECS**, WHO (**Enter CHEMLIST File for up-to-date regulatory information)

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ð IN FILE CA (1907 TO DATE)
TO NON-SPECIFIC DERIVATIVES IN FILE
IN FILE CAPLUS (1907 TO DATE)
IN FILE CAOLD (PRIOR TO 1967) 6224 REFERENCES 1 665 REFERENCES 1 6235 REFERENCES 1

2 REFERENCES

S Idarubicin/CN 1 IDARUBICIN/CN

=> D L10 1

LIO ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN RN 58957-92-9 REGISTRY
CN 5.12-Naphthacenedione, 9-acetyl-7-{(3-amino-2,3,6-trideoxy-α-L-1yxo-hexopyranosyl)cvy}]-7,8,9,10-tetrahydro-6,9,11-trihydroxy-, (7S,9S)- (9CI) OTHER CA INDEX NAME):

CN , 5,12-Naphthacenedione, 9-acety1-7-[(3-amino-2,3,6-trideoxy-α-L-lyxo-hexopyranosy1)oxy]-7,8,9,10-tetrahydro-6,9,11-trihydroxy-, (75-cis)-

OTHER NAMES:

1-Demethoxydaunorubicin 4-Demethoxydaunomycin 4-Demethoxydaunorubicin 4-DMD 22422222222

Demethoxydaunorubicin Idarubicin

STEREOSEARCH C26 H27 N 09 COM STN Files:

ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS,

BIOTECHNO, CA, CAPLUS, CASREACT, CENB, CHEMCATS, CIN, CSCHEM, DDFU, DRUGU, EMBAGE, IFICDB, ITIPAT, IFITUDB, IMSCOERANCH, INSDUGNEMS, INSPATENTS, IMSRESEARCH, IPA, WEDLINE, MRGK*, PATDRASPC, PHAR, PROMIT, PROUSDDR, PS, RIECS*, SYNTHLINE, TOXCENTER, USAN, USPATZ, USPATFULL,

(*File contains numerically searchable property data) Sources: WHO Other Sources:

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1391 REFERENCES IN FILE CA (1907 TO DATE) 87 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA 1394 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> S Labetalol/CN L11 1 LABETALOL/CN

-> D L11 1

ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 36894-69-6 REGISTRY SE ELI

Entered STN: 16 Nov 1984
Benzamide, 2-hydroxy-5-[1-hydroxy-2-[(1-methyl-3-phenylpropyl)amino]ethyl}(9CI) (CA INDEX NAME)

CON PACON DE CON PACON DE COND.

Albetol Ibidomide

Labetalol
C19 H24 N2 O3
C19 H24 N2 O3
STN Files: ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, COM
STN Files: ADISNEWS, CARRACT, CRNB. CHEMCATS, CHEMCATST, CIN. GSCHEM, DDFU,
CAA, CAPLUS, CASRACT, CRNB. FICTOB, IFIPAT, IFUUB, IMSPATENTS, IPA, MEDLINE,
MRCK*, NARALLERT, PRAR, PROWT, PS, RTEGS*, SPECINFO, SYNTHLINE,
TOXCENTER, USAN, USPATZ, USPATFULL
(*File contains numerically searchable property data)
Other Sources: EINEGS**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1007 REFERENCES IN FILE CA (1907 TO DATE)
26 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
1009 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> S Niclosamide/CN L12 1 NICLOSAMIDE/CN

=> D L12 1

ANSWER I OF I REGISTRY COPYRIGHT 2006 ACS on STN
CO-65-7 REGISTRY
Entered STN: 16 Nov 1984
Benzamide, 5-chloro-N-(2-chloro-4-nitrophenyl)-2-hydroxy- (9CI)

L12 CN ED

(CA INDEX CA INDEX NAMES: Salicylanilide, 2',5-dichloro-4'-nitro- (8CI) OTHER

12. 5-Dichloro-4'-nitrosalicylanilide
12-Chloro-4-nitrophenylamide-6-chlorosalicylic acid
2-Chloro-4-nitrophenylamide-6-chlorosalicylic acid
2-Hydroxy-5-Chloro-4'-nitrophenyl) benzamide
5-Chloro-2'-chloro-4'-nitrophenyl) salicylamide
5-Chloro-N-(2'-chloro-4'-nitrophenyl) salicylamide
5-Chloro-N-(2'-chloro-4'-nitrophenyl) salicylamide
BAY 2353
Bayer 2353
Bayluscid
Cestocide
Cestocide
Devermine
Fedal-relmin

Fenasal

Helmiantin HL 2447

Iomesan

Mansonil

Mato N-[2'-Chloro-4'-nitropheny]]-5-chlorosalicylamide N-(2-Chloro-4-nitrophenyl)-5-chlorosalicylamide

Niclocide

Niclosamide NSC 178296 Phenasal Radeverm

STN Files: ADISNEMS, AGRICOLA, ANABSTR, AQUIRE, BELLSTEIN*, BIOSIS, BIOTECHNO, CA, CABA, CAOLD, CAPLUS, CASREACT, CRNB, CHEMCATS, CHEMLIST, CIN, CSCHEM, CSNB, DDFU, DRUGU, EMBASE, HSDB*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, MSDS-OHS, PIRA, RROWT, PS, RFECS*, SCISEBACH, SPECINFO, TOXCENTER, TULSA, ULIDAT, USAN, USPATZ, USPATFULL, VETU (*File contains numerically searchable property data)
Other Sources: ENESC**, WHO

(**Enter CHEMLIST File for up-to-date regulatory information) Zestocarp 3D CONCORD 12687-52-4 C13 H8 C12 N2 O4 COM STN Files: ADISN BIOTECHNO, CA, C Sulqui Tredemine Utosamide Vermitid Vermitin WR 46234 Ruby Sagimid Yomesan

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

IN FILE CA (1907 TO DATE)
TO NOW-SPECIFIC DERIVATIVES IN FILE CA
IN FILE CAPLUS (1907 TO DATE)
IN FILE CAOLD (PRIOR TO 1967) 521 REFERENCES 1 12 REFERENCES 1 521 REFERENCES 1 13 REFERENCES 1

-> S Osalmid/CN 113 1 OSALMID/CN

-> D L13 1

L13 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN
RN 526-18-1 REGISTRY
ED Entered STN: 16 Nov 1984
CN Benzamide, 24-ydroxyphenyl)- (9CI) (CA INDEX NAME)
CN Salicylanilide, 4'-hydroxyphenyl) benzamide
CN 2-Hydroxyphenyl) benzamide
CN 2-Hydroxypalicylanilide
CN Auxobil
CN Bilone
CN Bilocol
CN Bilocol
CN Dribazil
CN Dribazil
CN Driol-Labaz
CN Endran

Fumispore Jestmin Kanochol L 1718

Li Dan Feng N'-Salicyloyl-p-aminophenol N-(4-Hydroxyphenyl)salicylamide N-(p-Hydroxyphenyl)salicylamide NSC 93960

Oksafenamide

Osalmid Osalmide

Oxaphenamid Oxaphenamide

p'-Hydroxysalicylanilide p-Hydroxyphenylsalicylamide

Salmidochol

STN Files: ANABSTR, BEILSTEIN*, BIOSIS, CA, CAOLD, CAPLUS, CASREACT, CHEMIST, CHEMIST, CIN, CSCHEM, DDFU, DRUGU, EMBASE, IFTCDB, IFTPAT, IFTUDB, IPA, MEDLINE, MECK*, RTECS*, SPECINFO, TOXCENTER, USAN, USPATFULL, VETU

(**TILe contains numerically searchable property data)
Other Sources: EINECS**, WHO

(**Enter CHEMIST File for up-to-date regulatory information) Saryuurin WR 17456 Yoshicol 3D CONCORD C13 H11 N 03 COM STN Files: A

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

IN FILE CA (1907 TO DATE)
IN FILE CAPLUS (1907 TO DATE)
IN FILE CAOLD (PRIOR TO 1967) 141 REFERENCES I 141 REFERENCES I 17 REFERENCES I

â

S Salacetamide/CN

-> D L14 1

L14 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 487-48-9 REGISTRY
CD Entered STN: 16 Nov 1984
CN Benzamde, N-acetyl-2-hydroxy- (9CI) (CA INDEX NAWE)
OTHER CA INDEX NAMES:
OTHER NAMES:
CN Zalicylamide, N-acetyle- (6CI, 7CI, 8CI)
OTHER NAMES:
CN Z-Hydroxy-N-acetylbenzamide
CN Acetsalicylamide
CN Acetsalicylamide
CN Acetsalicylamide

Salacetanide
Salacetanide
Salacetanide
Salacetanide
Salacyl
Salacyl
SD CONCORD
C9 H9 N O3

COM
STN Files: BELLSTEIN*, BIOSIS, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU, DRUGU, EMBASE, MRCK*, PROMT, PS, RTECS*, SPECINFO, TOXCENTER, USAN, USPATZ, USPATFULL

(**File contains numerically searchable property data)
Other Sources: EINEGS**, WHO

(**Enter CHEMLIST File for up-to-date regulatory information) Arthrisin Ethrisin Labazyl N-Acetylsalicylamide Nacemide NSC 40153 NSC 525079 NSC 631657 Rixamone

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

103 REFERENCES IN FILE CA (1907 TO DATE)
7 REFERENCES TO NON-SPECIFIC DEALTWATIVES IN FILE CA
103 REFERENCES IN FILE CAPLUS (1907 TO DATE)
8 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

L15 Salicylamide/CN
L15 1 SALICYLAMIDE/CN

-> D L15 1

ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 65-45-2 REGISTRY Entered STN: 16 Nov 1984 Benzamide, 2-hydroxy- (9CI) (CA INDEX NAME) R CA INDEX NAMES: 8.251cylamide (8CI)

2-Carbbamoylphenol 2-Carbbamoylphenol 2-Hydroxybenzamide Acket Acket Afro-Sal Algamon Algamon Algamida Allevin Amid-Sal Amidosal Amamid Benesal Cetamide

Dropsprin Liquiprin Cidal Cymidon

Morsarinas

Novecyl NSC 3115 NSC 83150 o-Hydroxybenzamide

Oramid Panithal Raspberin Salamid

Saliamid Saliamin Salicim Salicylic acid amide Salipur Salizell Salizell Salzin Salzin

Serramida SR 4326

Urtosal 3D CONCORD C7 H7 N O2

TH FILES: ADISNEMS, AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CRBA, CAOLD, CAPLUS, CASREACT, CBNB, CHEMATS, CHEMINFORMRX, CHEMINFORMRX, CHEMINFORMRX, CHEMINFORMRX, CHEMIST, CIN, CSCHEM, DDFU, DETHERM*, DRUGU, EMBASE, PRELIN*, HSDB*, IFICDB, IFIRT, IFIUDB, IPA, MEDLINE, MRCK*, PIRA, BROWT, PS, RTECS*, SPECINFO, SYNTHLINE, TOXCENTER, TULSA, USAN, USPATZ, VIB STN Files:

(*File contains numerically searchable property data)
Other Sources: DSL**, EINECS**, TSCA**, WHO
(**Enter CHEMLIST File for up-to-date regulatory information)

CI NH2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2122 REFERENCES IN FILE CA (1907 TO DATE)
77 REFERENCES IO NON-SPECIFIC DERIVATIVES IN FILE CA
7 REFERENCES IN FILE CAPLUS (1907 TO DATE)
7 REFERENCES IN FILE CADLD (FRIOR TO 1967)

s> S Salicylanilide/CN
L16
1 SALICYLANILIDE/CN

-> D L16

L16 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN

C13 H11 NO2

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CABLES ADISNEWS, AGRICOLA, ANABSTR, AQUIRE, BELLSTEIN*, BIOSIS, CA,

CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMEX, CHEMLIST, CIN, CSCHEM,

CSNB, DDFV, DETHERAY, DRUGU, BERBASE, GREINY, IFICOB, IFIPAT, IFIUDB,

IPA, MEDLINE, MRCK', MSS-OHS, PIRA, PROMT, RECS*, SPECINFO, SYNTHLINE,

TOXCENTER, USAN, USPATZ, USPATFULL, VETU, VTB

C*FILE contains numerically searchable property data)

Other Sources: DSL**, EINECS**, TSCA**

(**Enter CHEMLIST File for up-to-date regulatory information) (CA INDEX NAME) Entered STN: 16 Nov 1984
Benzamide, 2-hydroxy-N-phenyl- (9CI)
CA INDEX NAMES: 2-(N-Phenylcarboxamido) phenol 2-(Phenylaminocarbonyl) phenol 2-Hydroxy-N-phenylbenzamide 2-Hydroxybenzamilide Ansadol N-Phenyl-2-hydroxybenzamide N-Phenylsalicylamide NSC 14881 SA 88
Salicylanilid
Salicylic acid anilide
Salifebrin
Saliide Salicylanilide (8CI) o-Hydroxybenzanilide 87-17-2 REGISTRY Shirlan Shirlan AG Shirlan Extra WR 10019 3D CONCORD Salnide Sherstat SLN Aseptolan RN 97-17-2
CN Benzamid
OTHER CA INDE
OTHER NAMES:
CN 2-(N-Phe
CN 2-(Pheny)
CN 2-(Pheny)
CN 2-(Pheny)
CN 2-(Pheny)
CN 2-(Pheny)
CN 2-(Pheny)
CN Asadol
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PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

815 REFERENCES IN FILE CA (1907 TO DATE)
108 REFERENCES TO NON-SPECIFIC DEALYATIVES IN FILE CA
816 REFERENCES IN FILE CAPLUS (1907 TO DATE)
27 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> S Salsalate/CN Ll7 l SALSALATE/CN

-> D L17

Salsalate

N Salsalate

N Salsalate

N Salsalate

N Saspyrine

Sasspyrinum

Sasspyrinum

SURCORORD

STATE ADISNEWS, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CHEMLIST, CIN, CSCHEM, DDEU, DRUGH, EMBASE, ITICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, PROMT, PS, RTECS*, SCIESRACH, TOXCENTER, USAN, USAPATFULL

(*File contains numerically searchable property data)

Other Sources: EINECS**, NDSL**, TSCA**, WHO

(**Enter CHEMLIST File for up-to-date regulatory information) ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 552-94-3 REGISTRY Entered STN: 16 Nov 1984 Benzoic acid, 2-hydroxy-, 2-carboxyphenyl ester (9CI) (CA INDEX NAME) CA INDEX NAMES: Salicylic acid, bimol. ester (7CI) Salicylic acid, salicylate (6CI) Diplosal Disalgesic Disalgesic Disalicylic acid Disalicylic acid Mono-Gesic Nobacid NSC 49171 O-Salicylsalicylic acid Salicyl salicylate Salicyloxysalicylic acid Salicyloylsalicylic acid Salicylsalicylic acid Diacesal

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

375 REFERENCES IN FILE CA (1907 TO DATE)
21 REPRENCES TO NON-SPECIFIC DESTVATIVES IN FILE CA
21 REFERENCES IN FILE CAPIUS (1907 TO DATE)
311 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

-> S Zorubicin/CN 118 1 20RUBICIN/CN

=> D L18

Entered STN: 16 Nov 1984
Bensofc acid, [1-[(25,48)-4-[(3-amino-2,3,6-trideoxy-a-L-1yxo-dioxy-apyranosyl)oxy]-1,2,3,4,6,11-hexahydro-2,5,12-trihydroxy-7-methoxy-6,11-dioxo-2-naphthacenyllethylidene]hydrazide (9CI) (CA INDEX NAME) Benzoic acid, [1-[4-[(3-amino-2,3,6-trideoxy-u-L-1yxo-hexopyranosyl)oxy]-1,2,3,4,6,11-hexahydro-2,5,12-trihydroxy-7-methoxy-6,11-dioxo-2-naphthacenyl]ethylldene]hydrazide, (25-cis)-TN Files: ADISNEWS, ANABSTR, BELLSTEIN*, BIOSIS, BIOTECHNO, CA, CAPLUS, CHENCATS, CIN, DDFU, DBUGU, EMBASE, INSPATENTS, IRA, MEDLINE, MRCK*, MRSALLERT, PHAR, PROMT, PROUSDDR, PS, RTECS*, SYNTHLINE, TOXCENTER, USAN, USPAT, USPATFULL.

(*File contains numerically searchable property data) REGISTRY COPYRIGHT 2006 ACS on STN Daunomycin benzoylhydrazone C34 H35 N3 O10 COM STN Files: AD Other Sources: ANSWER 1 OF 1 54083-22-6 RE Zorubicin STEREOSEARCH Rubidazone MCMC 4222 Rubidazon NAMES: OTHER CN B 2 2 2 3 3 2 3 3 3

(CA INDEX NAME)

9,10-Anthracenedione, 1,2-dihydroxy- (9CI)

CA INDEX NAMES

Absolute stereochemistry. Double bond geometry unknown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

165 REFERENCES IN FILE CA (1907 TO DATE) 8 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA 168 REFERENCES IN FILE CAPLUS (1907 TO DATE)

Other Sources:

=> S Alizarin/CN L19 1 ALIZARIN/CN

-> D L19 1

ANSWER I OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 72-48-0 REGISTRY Entered STN: 16 Nov 1984 L19 RN ED

1,2-Dibydroxy-9,10-anthracenedione
1,2-Dibydroxy-9,10-anthraquinone
1,2-Dibydroxy-9,10-anthraquinone
1,2-Dibydroxy-9,10-anthraquinone
1,2-Dibydroxy-9,10-anthraquinone
1,2-Dibydroxy-9,10-anthraquinone
1,2-Dibydroxy-9,10-anthraquinone
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1,2-Dibydroxy-9,10-anthraquinone
1,2-Dibydroxy-9,10-anthraquinon Anthraquinone, 1,2-dihydroxy- (BCI) 1,2-Anthraquinonediol Alizarin B (6CI STN Files:

REFERENCES IN FILE CA (1907 TO DATE)
REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
REFERENCES IN FILE CAPLUS (1907 TO DATE)
REFERENCES IN FILE CAOLD (PRIOR TO 1967) 1798 129 1804 15

S Alkannin/CN 2 ALKANNIN/CN

■> D L20 1-2

ANSWER 1 OF 2 REGISTRY COPYRIGHT 2006 ACS on STN
2344-65-7 REGISTRY
Entered STN: 16 Nov 1984
1,4-Naphthalenedione, 5,8-dihydroxy-6-[(1S)-1-hydroxy-4-methyl-3-pentenyl]-(CA INDEX NAME)

1,4-Naphthalenedione, 5,8-dihydroxy-6-(1-hydroxy-4-methyl-3-pentenyl)-, OTHER CA INDEX NAMES CN 1,4-Naphthaleneo

1,4-Naphthoquinone, 5,8-dihydroxy-6-(1-hydroxy-4-methyl-3-pentenyl)-, (-)-S

Alkanet extract Alkanet extract

extract, inspissated

Alkanna red

Alkannin

Anchusa acid

Anchusin

C.I. 75530 C.I. Natural Red 20 STEREOSEARCH

C16 H16 05
STN FILES: AGRICOLA, BELLSTEIN*, BIOSIS, BIOTECHNO, CA, CAPLUS, CHEMCATS, CHEMLIST, EMBASE, MEDLINE, RTECS*, TOXCENTER, USPATFULL (*File contains numerically searchable property data)

Absolute stereochemistry

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

17 REFERENCES IN FILE CA (1907 TO DATE) 17 REFERENCES IN FILE CAPLUS (1907 TO DATE)

ANSWER 2 OF 2 REGISTRY COPYRIGHT 2006 ACS on STN 517-88-4 REGISTRY Entered STN: 16 Nov 1984 E S E

1,4-Naphthalenedione, 5,8-dihydroxy-2-[(18)-1-hydroxy-4-methyl-3-pentenyl]-(9CI) (CA INDEX NAME) [6]1,4-Naphthoquinone, 5,8-dihydroxy-2-(1-hydroxy-4-methyl-3-pentenyl)-, (-)-STN Files: ACRICOLA, ANABSTR, BEILSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMRX, CHEMLIST, CSCHEM, DDFU, DRUGU, ENBASE, IPA, MRCK*, SPECINFO, TOXCENTER, USPATFULL (*File contains numerically searchable property data)
Other Sources: EINECS**
(**Enter CHEMLIST File for up-to-date regulatory information) OTHER CA INDEX NAMES: CN 1,4-Naphthalenedione, 5,8-dihydroxy-2-(1-hydroxy-4-methyl-3-pentenyl)-, 20 Alkanet extract Alkanna red Anchusa acid Anchusin C.1. 7530 C.1. Natural Red 20 NSC 94524 (8CI) Alkannin (6CI) STEREOSEARCH)-Alkannin C16 H16 O5 CON SECONDARY SE S

Absolute stereochemistry. Rotation (-).

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

145 REFERENCES IN FILE CA (1907 TO DATE)
16 REPERSENCES TO NON-SPECTEIC DEFENTATIVES IN FILE CA
17 REFERENCES IN FILE CAPLUS (1907 TO DATE)
17 REFERENCES IN FILE CADLO (PRIOR TO 1967)

S Anthragallol/CN 1 ANTHRAGALLOL/CN

-> D L21 1

121 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN
RN 602-64-2 REGISTRY
ED for 602-64-2 REGISTRY
CN 9,10-Anthracenedione, 1,2,3-trihydroxy- (9CI) (CA INDEX NAME)
OTHER CA INDEX NAMES:
CN Anthragallol (6CI)
CN Anthragallone, 1,2,3-trihydroxy- (7CI, 8CI)
OTHER NAMES:
CN 1,2,3-Trihydroxyanthraquinone
CN 1,2,3-Trihydroxyanthraquinone
CN 1,2,3-Trihydroxyanthraquinone
CN Alizarine Brown HD

STN Files: AGRICOLA, ANABSTR, BELLSTEIN*, BIOSIS, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CSCHEW, MRCK*, NAPRALERT, RTECS*, SPECINFO, TOXCENTER, USPATFULL.

(*File contains numerically searchable property data)
Other Sources: EINECS**

(**Enter CHEMLIST File for up-to-date regulatory information) Alizarine Brown R
Anthracene Brown FD
Anthracene Brown G
Anthracene Brown G
Anthracene Brown N
Anthracene Brown WH
Anthracene Brown WH
Anthracene Brown WL
Anthracene Printing Brown
Antracene Printing Brown
Antragallol
C.I. \$8200
C.I. Mordant Brown 42
C.I. Mordant Brown 42
C.I. Mordant Brown A2
C.I. Mordant Brown A3
CONCORD
C.I. Mordant Brown GC
Mitsul Anthracene Brown
NSC 31754
3D CONCORD
CI4 H8 05
COM
STN Files: AGRICOLA, ANABS

32073-06-6 C14 H8 O4 COM STN Files:

문 A C C

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

REFERENCES IN FILE CA (1907 TO DATE)
REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA REFERENCES IN FILE CAPLUS (1907 TO DATE)
REFERENCES IN FILE CADLD (PRIOR TO 1967) 93 21 21

=> S Anthrarufin/CN L22 1 ANTHRARUFIN/CN

-> D L22 1

(CA INDEX NAME) L22 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS ON STN
RN 117-12-4 REGISTRY
ED FILECAGE STN: 16 Nov 1984
CN 9,10-Anthracenedione, 1,5-dihydroxy- (9CI) (CA INDEX OTHER CA INDEX NAMES:
CN Anthraquinone, 1,5-dihydroxy- (8CI)
CN Anthraufin (6CI)
OTHER NAMES:
CN 1,5-Dihydroxy-9,10-anthraquinone
CN 1,5-Dihydroxy-9,10-anthraquinone
CN 1,5-Dihydroxyanthraquinone
CN NSC 646570
CN NSC 646570
CN NSC 646570
CN NSC 646570
CN NSC 646570
CN NSC 646570

Laxanthreen Laxipur Laxipurin Laxanorm

Dorbane Istizin

STN Files: AGRICOLA, ANABSTR, AQUIRE, BEILSTEIN*, BIOSIS, CA, CAOLD, CARLES, CARALTATS, CHEMATAY, CHEMILTST, CSCHEM, DDEU, DETHERN*, DRUGU, GMELIN*, HSDB*, IFTCB, IFTCB, IFTUB, HSDF, MSDS-OHS, NAPRALERT, RIECS*, SPECINFO, TOXCENTER, USPATZ, USPATFULL ('File contains numerically searchable property data) Other Sources: EINEGS**, NDSL**, TSCA** (**Enter CHEMIST File for up-to-date regulatory information) ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN
117-10-2 REGISTRY
ED Entered STN: 16 Nov 1984
117-10-2 REGISTRY
The CA INDEX NAMES:

NATH ANTHROUGHORD, 1,8-dihydroxy- (9CI) (CA INDEX NAME)
THER CA INDEX NAMES:

NATHER NAMES:
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NAMES: 466 REFERENCES IN FILE CA (1907 TO DATE)
13 REFERENCES TO NON-SPECIFIC DEFIVATIVES IN FILE CA
466 REFERENCES IN FILE CADLUS (1907 TO DATE)
27 REFERENCES IN FILE CADLUS (PRIOR TO 1967) **PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT** => S Danthron/CN L23 1 DANTHRON/CN => D L23 1

Switsalax
32073-00-7,
343235-40-5
C14 H8 04
C17 H8 04
C18 H8 04
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C1 Modane NSC 38626 NSC 646568 NSC 7210

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

985 REFERENCES IN FILE CA (1907 TO DATE)
54 REFERENCES TO NON-SPECIFIC DERLYATIVES IN FILE CA
985 REFERENCES IN FILE CAPLUS (1907 TO DATE)
2 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> S emodin/CN L24 2 EMODIN/CN -> D L24 1-2 ANSWER 1 OF 2 REGISTRY COPYRIGHT 2006 ACS on STN
LS697-21. REGISTRY
Entered STN: 16 Nov 1984
Benzeneacetic acid, \alpha-methyl-4-(2-methylpropyl)- (9CI) (CA INDEX S 5 5 5 5

(#)-d-Methyl-4-(2-methylpropyl)benzeneacetic acid (#)-1buprofen (#)-lbuprophen (#)-lbuprophen (4-Isobutylphenyl)-d-methylacetic acid (RS)-Ibuprofen OTHER NAMES:

CN (1)-α-Met
CN (1)-1-bup
CN (1)-1bup
CN (4)-1bup
CN (4)-1sobut
CN (4-1sob
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a-(4-isoburylphenyl)propionic acid a-Methyl-4-(2-methylpropyl)benzeneacetic acid 2-(4'-isoburylphenyl)propionic acid 2-(4-isoburylphenyl)propanoic acid 2-(p-isoburylphenyl)propanoic acid 4-isoburylphenyl)propionic acid 4-isoburylhydratropic acid A-isoburylhydratropic acid

Adex 200

BIOTECHNO. CA. CABA, CANOLD, CARREACT. CBNB, CEMECATS,
BIOTECHNO. CA. CABA, CANOLD, CAPLUS, CARREACT. CBNB, CEMECATS,
CHEMINEDRAKX, CHEMIST, CIN, CSCHEM, CSNB, DDTU, DRUGU, EMBASE, HSDB.,
IFICOB, IFIPAT, IFTUDB, IMSCOSEARCH, IMSDRUGNEMS, IMSPATENTS,
IMSESERACH, IRP, MEDLINE, MRCK*, MSDS-OHS, PATDPRASC, PHAR, PIRA,
PROMT, PROUSDDR, PS, RECS*, SCISSARCH, SPECINFO, SYNTHLINE, TOXCENTER,
ULIDAT, USAN, USPAT2, USPATFULL, VETU

(**Elle contains numerically searchable property data)
Other Sources: DSL**, Elle for up-to-date regulatory information) CN Adran
CN Advin
CN Advin
CN Algofen
CN American
CN Amican
CN Antifen
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g 8613 REFERENCES IN FILE CA (1907 TO DATE)
249 REFERENCES TO NON-SPECIFIC DERLYAPITIES IN FILE
8634 REFERENCES IN FILE CAPLUS (1907 TO DATE)
2 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

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CLS HIO OS
CONCORD
STATE FILES:
ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BELLSTEIN*, BIOSIS,
BIOTECHNO, CA, CABA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMINFORMEX,
CHEMIJST, CIN, CSCHEM, CSNSB, DDFU, DETHERM*, DRUGU, EMBASE, HSDB*,
IMSCOSEARCH, IMSDRUGNEWS, IMSRESEARCH, IPA, MEDLINE, MRCK*, NAPRALERT,
PROMIT, PROUSDBR, RTECS*, SPECINFO, TOXCENTER, USPATZ, USPATFULL
(*File contains numerically searchable property data)
Other Sources: EINECS**
(**Enter CHEMIST File for up-to-date regulatory information)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1647 REFERENCES IN FILE CA (1907 TO DATE)
51 REFERENCES TO NON-SPECIFIC DEMINATIVATIVES IN FILE CA
1649 REFERENCES IN FILE CAPLUS (1907 TO DATE)
23 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> S Purpurin/CN

2 PURPURIN/CN

=> D 125 1-2

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STN Files: AGRICOLA, ANABSTR, AQUIRE, BELLSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CARLUS, CARREACT, CHEMATS, CHEMISTY, CHIN, CSCHEM, DDFU, DETHERH*, DRUGU, EMBASE, GWELIN*, IFICDB, IFIPAT, IFIUDB, IPA, MEDLINE, MRCK*, PIRA, PROMT, RTECS*, SPECINFO, TOXCENTER, USPATZ, USPATZLL (*File contains numerically searchable property data) other Sources:
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                             N Files: AGRICOLA, BIOSIS, CA, CAPLUS, CHEMCATS, CIN, IFICDB, IFIUDB, PIRA, PROMT, TOXCENTER, USPATZ, USPATFULL
L25 ANSWER 1 OF 2 REGISTRY

RN 75775-33-6 REGISTRY

ED Entered STN: 16 Nov 1984

CN 4H-Fuzo(3', 2'.4,5)fuzo(2,3-h)-1-benzopyran-4-one, 10-(acetyloxy)-
2.3,73,9,10,10a-hexahydro-9,9-dimethyl-2-phenyl- (9CI) (CA INDEX NAME)

OTHER NAMES:
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                    LUSS ANSWER 2 OF 2 REGISTRY COPYRIGHT 2006 ACS on STN
RN 81-54-9 REGISTRY
ED Entered STN: 16 Nov 1984
CN 9.10-Anthreachedone, 1,2,4-trihydroxy- (9CI) (CA INDEX NAWE)
CN Anthraquinone, 1,2,4-trihydroxy- (7CI, 8CI)
CN Anthraquinone, 1,2,4-trihydroxy- (7CI, 8CI)
CN 1,2,4-Trihydroxy-9,10-anthraquinone
CN 1,2,4-Trihydroxy-9,10-anthraquinone
CN 1,2,4-Trihydroxy-9,10-anthraquinone
CN 1,2,4-Trihydroxy-9,10-anthraquinone
CN 1,2,4-Trihydroxy-9,10-anthraquinone
CN 1,2,4-Trihydroxy-9,10-anthraquinone
CN 1,2,4-Trihydroxy-9,10-anthraquinone
CN 1,2,4-Trihydroxy-9,10-anthraquinone
CN 1,2,4-Trihydroxy-9,10-anthraquinone
CN 1,2,4-Trihydroxy-9,10-anthraquinone
CN 1,2,4-Trihydroxy-9,10-anthraquinone
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CN 1,2,4-Trihydroxy-1,2,10-anthraquinone
CN 1,2,4-Trihydroxy-1,2
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18 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE
37 REFERENCES IN FILE CAPLUS (1907 TO DATE)
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                        **PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT**
                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                                            Currently available stereo shown
                                                                                                                                                                                                                                                                                                                                                                                                                             Purpurin (flavanone)
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C23 H22 O6
STN Files: P
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STN Files: AGRICOLA, ANABSTR, AQUIRE, BELLSTEIN*, BIOSIS, BIOTECHNO, CA, CAOLD, CARLUS, CASREACT, CHEMACATS, CERMILST, CSCHEM, DBTU, DBRUU, ENGASE, GMELLIN*, IFICDB, IFIPAT, IFIUDB, IPA, MRCK*, MSDS-OHS, NAPRALERT, RIECS*, SPECINFO, TOXCENTER, USPATZ, USPATFULL (*File contains numerically searchable property data) other Sources: DSL**, EINECS**, TSCA** (**Enter CHEMLIST File for up-to-date regulatory information)
                                                                                                                                                                                                                                          ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN 81-61-8 REGISTRY
Entered STN: 16 Nov 1984
9,10-Anthracenedione, 1,2,5,8-tetrahydroxy- (9CI) (CA INDEX NAME):
A CA INDEX NAMES:
A Anthraquinone, 1,2,5,8-tetrahydroxy- (8CI)
Quinalizarin (6CI)
537 REFERENCES IN FILE CA (1907 TO DATE)
53 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
537 REFERENCES IN FILE CALLUS (1907 TO DATE)
48 REFERENCES IN FILE CADLU (PRIOR TO 1967)
                                                                                                                                                                                                                                                                                                                                                                                                             HEK NAMES.

1, 2, 5, 8-Tetrahydroxy-9, 10-anthraquinone

1, 2, 5, 8-Tetrahydroxyanthraquinone

1, 4, 5, 6-Tetrahydroxyanthraquinone

1, 4, 5, 6-Tetrahydroxyanthraquinone

1, 4, 5, 6-Tetrahydroxyanthraquinone

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Alizarine

NSC 144046

NSC 14066

Quinalizarine

3D CONCORD

COM

COM

STN Files: AGRICOLA, ANABSTR, AQUIRE, BE.

STN Files: AGRICOLA, ANABSTR, AQUIRE, BE.
                                                                                                                                     S Quinalizarin/CN
L26 1 QUINALIZARIN/CN
                                                                                                                                                                                                    => D L26 1
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PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

465 REFERENCES IN FILE CA (1907 TO DATE)
51 REPERENCES TO NON-SPECTIC DERIVATIVES IN FILE CA
465 REFERENCES IN FILE CAPLUS (1907 TO DATE)
30 REFERENCES IN FILE CADLD (PRIOR TO 1967)

1 QUINIZARIN/CN -> S Quinizarin/CN

=> D L27 1

81-64-1 REGISTRY
Entered STN: 16 Nov 1984
9,10-Anthracenedione, 1,4-dihydroxy- (9CI) (CA INDEX NAME)
CA INDEX NAMES: ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN Anthraquinone, 1,4-dihydroxy- (8CI) 127 ANSWER 1

EN 81-64-1

ED CO 9,10-ANT

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COM STN Files:

STUTION THES: ACRICOLA, ANABSTR, AQUIRE, BELLSTEIN*, BIOSIS, BIOTECHNO, CA, CADLUS, CASERACT, CHEMCATS, CHEMINFONRAX, CHEMLIST, CIN, CSCHEM, CSNB, DDEV, DETHERA*, DRUGU, EMBASE, GENBANK, GMELIN*, HSDB*, IFICDB, IFIPAT, IFIUDB, MEDLINE, MSC*, NSDS-OHS, NAPPALERT, PIRA, PROMT, RTECS*, SPECINFO, TOXCENTER, USPATFULL
(***IFI contains numerically searchable property data)
Other Sources: DSL**, FINECS**, TSCA**
(***Enter CHEMLIST File for up-to-date regulatory information)

1483 REFERENCES IN FILE CA (1907 TO DATE)
89 REFERENCES TO ONG-SPECIFIC DERIVATIVES IN FILE CA
1483 REFERENCES IN FILE CAPLUS (1907 TO DATE)
6 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> S "(Z)-3-hydroxyacrylaldehyde"/CN 128 0 "(Z)-3-HYDROXYACRYLALDEHYDE"/CN

-> S "3-hydroxyacrylaldehyde"/CN L29 0 "3-HYDROXYACRYLALDEHYDE"/CN

=> S hydroxyacrylaldehyde/CN
L30 0 HYDROXYACRYLALDEHYDE/CN

=> S "3-hydroxy-2-propene aldehyde"/CN 0 "3-HYDROXY-2-PROPENE ALDEHYDE"/CN

S hydroxypropenealdehyde/CN L32 0 HYDROXYPROPENEALDEHYDE/CN

=> S "3-hydroxy-2-propene-a1"/CN L33 0 "3-HYDROXY-2-PROPENE-AL"/CN